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İ		548/473.ccls. or 588/417.ccls. or		
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3	BRS	<b>L</b> 7	256	549/69.ccls.	USPAT; US-PG PUB; EPO; JPO; DERWE NT	2004/07/16 06:05			
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SINCE FILE TOTAL ENTRY SESSION 0.21

0.21

FULL ESTIMATED COST

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14 JUL 2004 HIGHEST RN 710278-45-8 STRUCTURE FILE UPDATES: DICTIONARY FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8

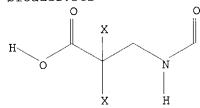
TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

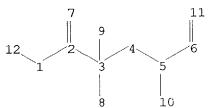
Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

Uploading C:\Examination Auxillary files\09995987\09995987 third try fixed H broader.str





chain nodes : 1 2 3 4 5 6 7 8 9 10 11 12 chain bonds : 3-9 4-5 5-6 5-10 6-11 1-2 1-12 2-3 2-7 3-4 3-8 exact/norm bonds : 4-5 5-6 6-11 exact bonds : 1-12 2-3 3-4 3-8 3-9 5-10 normalized bonds :

G1:C,O,S,N

1-2 2-7

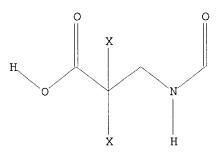
Hydrogen count : 4:>= minimum 2 Match level: 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS

## STRUCTURE UPLOADED L1

=> d 11

L1 HAS NO ANSWERS

STR L1



G1 C, O, S, N

Structure attributes must be viewed using STN Express query preparation.

=> search 11 sss sam SAMPLE SEARCH INITIATED 06:15:22 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED

21 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:

ONLINE \*\*COMPLETE\*\*

\*\*COMPLETE\*\* BATCH

PROJECTED ITERATIONS:

146 TO 694

PROJECTED ANSWERS:

1 TO 80

=> d scan

L2

REGISTRY COPYRIGHT 2004 ACS on STN 1 ANSWERS L2

1 SEA SSS SAM L1

Propanoic acid, 3-[[4-[[[(3,5-dichlorophenyl)amino]carbonyl]][4-[4-(1,1-dichlorophenyl)amino]carbonyl]]IN dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2difluoro- (9CI)

C34 H35 Cl2 F2 N3 O4 MF

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=> search l1 sss full FULL SEARCH INITIATED 06:15:46 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 525 TO ITERATE

100.0% PROCESSED 525 ITERATIONS SEARCH TIME: 00.00.01

3 ANSWERS

L3

3 SEA SSS FUL L1

=> d scan

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 3-[[4-[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro-(9CI)

MF C34 H35 C12 F2 N3 O4

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]- (9CI)

MF C11 H11 F2 N O4

$$\begin{array}{c} \text{O} \\ \parallel \\ \text{Ph-CH}_2 - \text{O-C-NH-CH}_2 - \text{CF}_2 - \text{CO}_2 \text{H} \end{array}$$

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 3-[[4-[[[4-(1-cyclohexen-1-yl)phenyl][[(3,5-

dichlorophenyl) amino] carbonyl] amino] methyl] benzoyl] amino] -2,2-difluoro-(9CI)

MF C30 H27 Cl2 F2 N3 O4

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL SESSION 155.84 156.05

FULL ESTIMATED COST

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FILE COVERS 1907 - 16 Jul 2004 VOL 141 ISS 4 FILE LAST UPDATED: 15 Jul 2004 (20040715/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 13

L4 1 L3

=> d 14 ti fbib abs

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN

TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.

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2002:391685 CAPLUS
ΝA
     136:385945
DN
      Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon
TI
      antagonists/inverse agonists.
      Jorgensen, Anker Steen; Madsen, Peter
IN
     Novo Nordisk A/S, Den.
PA
     PCT Int. Appl., 85 pp.
SO
      CODEN: PIXXD2
DT
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      English
LA
FAN.CNT 1
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                         KIND DATE
      PATENT NO.
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      ______
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      WO 2002040446
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RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG DK 2000-1733 A 20001117
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OS
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AΒ
      heteroarylene; X = (CH2)q(CR12R13)r(CH2)s, CO(CR12R13)r(CH2)s,
      NR11CO(CR12R13)r(CH2)s, etc.; r = 0, 1; s = 0-3; R11, R12, R13 = H,
      alkyl; D = (substituted) Ph, naphthyl, pyridyl, indenyl, benzothienyl,
      thienyl, furyl, benzofuryl, etc.; E = (substituted) cyclohexyl, Ph, PhCH2,
      PhCH2CH2, indanyl, benzhydryl, etc.], were prepared Thus, Me
      4-[(4-cyclohex-1-enylphenylamino)methyl]benzoate (preparation given) in CH2Cl2
      containing diisopropylethylamine was treated with 3,5-dichlorophenyl
      isocyanate to give a residue which was saponified with LiOH. The resulting
      acid in DMF was treated with 3-[(dimethyliminium)(dimethylamino)methyl]-
      1,2,3-benzotriazol-1-ium-1-olate hexafluorophosphate,
      diisopropylethylamine, Me 3-amino-2,2-difluoropropionate hydrochloride to
      give the uncharacterized amide ester, which was saponified with aqueous LiOH in
      THF/MeOH to give 3-[4-[1-(4-cyclohex-1-enylphenyl)-3-(3,5-
      dichlorophenyl)ureidomethyl]benzoylamino]-2,2-difluoropropionic acid. In
      a human glucagon receptor binding assay, title compds. showed IC50<1000
      nM.
                 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
RE.CNT 3
                 ALL CITATIONS AVAILABLE IN THE RE FORMAT
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FULL ESTIMATED COST 3.14 159.19

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=> file caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

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SINCE FILE ENTRY S 0.46

TOTAL SESSION 0.67

FULL ESTIMATED COST

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STRUCTURE FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8 DICTIONARY FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

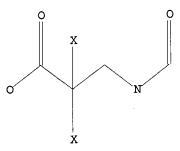
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> Uploading C:\Examination Auxillary files\09995987\09995987 patetable core broadest.str

L1 STRUCTURE UPLOADED

=> d l1 L1 HAS NO ANSWERS L1 STR



G1 C, O, S, N

Structure attributes must be viewed using STN Express query preparation.

=> search 11 sss sam
SAMPLE SEARCH INITIATED 07:24:18 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 194 TO ITERATE

100.0% PROCESSED 194 ITERATIONS

2 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS:

3045 TO 4715

PROJECTED ANSWERS:

2 TO 124

 $L_2$ 

2 SEA SSS SAM L1

=> d scan

L2 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)

MF C34 H35 Cl2 F2 N3 O4

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L2 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]-, methyl
 ester (9CI)

MF C12 H13 F2 N O4

$$\begin{array}{c} {\rm O} & {\rm O} \\ \parallel & \parallel \\ {\rm MeO-C-CF_2-CH_2-NH-C-O-CH_2-Ph} \end{array}$$

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=> search 11 sss full FULL SEARCH INITIATED 07:24:40 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 4048 TO ITERATE

100.0% PROCESSED 4048 ITERATIONS SEARCH TIME: 00.00.01

4 ANSWERS

\_\_\_

4 SEA SSS FUL L1

=> d scan

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]-, methyl
 ester (9CI)

MF C12 H13 F2 N O4

$$\begin{array}{c} {\rm O} & {\rm O} \\ \parallel & \parallel \\ {\rm MeO-C-CF_2-CH_2-NH-C-O-CH_2-Ph} \end{array}$$

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):4

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]- (9CI)

MF C11 H11 F2 N O4

$$\begin{array}{c} \text{O} \\ || \\ \text{Ph-CH}_2 - \text{O-C-NH-CH}_2 - \text{CF}_2 - \text{CO}_2 \text{H} \end{array}$$

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 3-[[4-[[[4-(1-cyclohexen-1-yl)phenyl][[(3,5-dichlorophenyl)amino]carbonyl]amino]methyl]benzoyl]amino]-2,2-difluoro-(9CI)

MF C30 H27 Cl2 F2 N3 O4

## \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)

MF C34 H35 Cl2 F2 N3 O4

## \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

## ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 155.84 156.51

FULL ESTIMATED COST

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FILE COVERS 1907 - 16 Jul 2004 VOL 141 ISS 4 FILE LAST UPDATED: 15 Jul 2004 (20040715/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

```
=> 13
            1 L3
L4
=> d 14
    ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
L4
    2002:391685 CAPLUS
AN
     136:385945
DN
     Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon
TI
     antagonists/inverse agonists.
     Jorgensen, Anker Steen; Madsen, Peter
ΙN
    Novo Nordisk A/S, Den.
PA
     PCT Int. Appl., 85 pp.
SO
     CODEN: PIXXD2
DT
     Patent
    English
LA
FAN.CNT 1
                                         APPLICATION NO. DATE
                     KIND DATE
     PATENT NO.
     _____
                                          ______
                                        WO 2001-DK760 20011115
     WO 2002040446 A1 20020523
PΙ
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             CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
             GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL,
             PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG,
             UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
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             CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR,
             BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
                      A5
                           20020527
                                         AU 2002-23502
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                           20030924
     EP 1345891
                      Α1
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             IE, SI, LT, LV, FI, RO, MK, CY, AL, TR
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     JP 2004513936
                     T2
                           20040513
                                          US 2001-995987
                                                           20011116
     US 2003027849
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                           20030206
PRAI DK 2000-1733
                      Α
                           20001117
     US 2000-252322P
                       Ρ
                           20001120
     WO 2001-DK760
                      W
                           20011115
     MARPAT 136:385945
              THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
RE.CNT 3
              ALL CITATIONS AVAILABLE IN THE RE FORMAT
```

=> file beilstein
COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
2.03 158.54

FILE 'BEILSTEIN' ENTERED AT 07:26:00 ON 16 JUL 2004 COPYRIGHT (c) 2004 Beilstein-Institut zur Foerderung der Chemischen Wissenschaften licensed to Beilstein GmbH and MDL Information Systems GmbH FILE RELOADED ON OCTOBER 20, 2002 FILE LAST UPDATED ON JUNE 15, 2004

FILE COVERS 1771 TO 2003.
\*\*\* FILE CONTAINS 8,997,153 SUBSTANCES \*\*\*

>>> PLEASE NOTE: Reaction data and substance data are stored in separate documents and can not be searched together in one query. Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a molecular formula or a structure search for example can be restricted to compounds with available reaction information by concatenation with PRE/FA, REA/FA or more general with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be selected from substance answer sets and searched in the next step as reaction partner BRNs - Reactant (RX.RBRN) or Product BRN (RX.PBRN). After a search for reaction details substance documents associated with reactants or products may be retrieved by searching RX.PBRNs or RX.RBRNs as BRNs. <<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

\*\*\*\*\*\*\*\*\*\*\*\*\*\*\*\*\*

\* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST.

\* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE

\* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE

\* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS.

0 ANSWERS

=> 11 SAMPLE SEARCH INITIATED 07:26:08 FILE 'BEILSTEIN' SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

100.0% PROCESSED 67 ITERATIONS SEARCH TIME: 00.00.02

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*
BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 849 TO 1831 PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L1